92197 SEARCH REQUEST FORM

Access DB# \_\_\_\_\_

# Scientific and Technical Information Center

Art Unit: 1624 Phor Mail Box and Bldg/Room Loca	ne Number 30 <u>8 4777</u> tion: <u>JD/5</u> R	Examiner #: 59/93 Date: 7  Serial Number: 09/93 esults Format Preferred (circle): PAPER Date: 100 pt.	1/23 C/49 DISK E-MAIL
Include the elected species or structure	es, keywords, synonyms, acr	be as specifically as possible the subject matter to be as specifically as possible the subject matter to be conyms, and registry numbers, and combine with the meaning. Give examples or relevant citations, autured abstract.	he concent or
Title of Invention:			
Inventors (please provide full names	s):		
	"· <del></del>		
Earliest Priority Filing Date:			<u>-</u>
	<u> </u>	<del></del>	
appropriate serial number.	O hord	p (parent, child, divisional, or issued patent numbers)	along with the
C	- G / what	HIC	
N .		N Con	1-A
0	N N	<i>K</i> /	
Tech.	Mary Jane Ruhl Info. Specialist, STIC TC-1600 W-1, Room 6A-06 hone: 605-1155	A=H/cts No py 2002 a 2	2003
	******		*****
STAFF USE ONLY	Type of Search	Vendors and cost where applicable	
earcher:	NA Sequence (#)	STN	<u>·</u>
	AA Sequence (#)		<del></del>
earcher Location:		Questel/Orbit	•
Pate Searcher Picked Up:			
	Litigation	Lexis/Nexis	<del></del>
earcher Prep & Review Time:	<u> </u>	Sequence Systems	<del></del>
Iterical Prep Time:	Patent Family	WWW/Internet	

Other (specify)

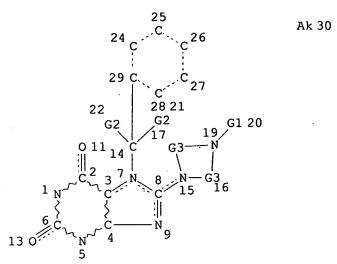
PTO-1590 (8-01)

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FILE 'REGISTRY' ENTERED AT 10:12:32 ON 24 APR 2003
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L1
L2
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L3
                   STR L3
1.4
L5
                   STR L4, DIS
L6
                   STR L4
                   STR L6
L7
                 0 S L7
L8
L9
                 O S L7 FULL
L10
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                 0 S L10
L11
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L12
L13
                   STR L10
L14
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L15
L16
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L18
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                   STR L19
L20
     2 S L20 FULL 2 compds in Reg. -pls see done stat for structure
FILE 'HCAPLUS' ENTERED AT 10:56:37 ON 24 APR 2003
1 S L22 / cit in CA Plus - its 2002, attacked
                 1 S L20
L21
L22
L23
                         (Ocits prior to 2002)
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=> d que stat 123 L20 STR



VAR G1=H/CH3
VAR G2=C/H
REP G3=(1-7) CH2
NODE ATTRIBUTES:
DEFAULT MLEVEL IS ATOM
GGCAT IS LOC UNS AT 30
DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES: RING(S) ARE ISOLATED OR EMBEDDED NUMBER OF NODES IS 26

STEREO ATTRIBUTES: NONE

L22 2 SEA FILE=REGISTRY SSS FUL L20 L23 1 SEA FILE=HCAPLUS ABB=ON L22

### => d ibib abs hitstr 1-1

L23 ANSWER 1 OF 1 HCAPLUS COPYRIGHT 2003 ACS ACCESSION NUMBER: 2002:31449 HCAPLUS

DOCUMENT NUMBER: 136:102402

TITLE: Preparation of piperazinylpurinediones as inhibitors

of dipeptidylpeptidase IV.

INVENTOR(S): Kanstrup, Anders Bendtz; Christiansen, Lise Brown;

Lundbeck, Jane Marie; Sams, Christian K.; Kristiansen,

Marit

PATENT ASSIGNEE(S): Novo Nordisk A/S, Den.

SOURCE: PCT Int. Appl., 135 pp.

CODEN: PIXXD2
DOCUMENT TYPE: Patent

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

P	PATENT NO.			KI	ND DATE				APPLICATION NO.						DATE			
W(	WO 2002002560				 2	2002	0110		1	 WO 20	01-D	 K467		2001	0704			
	W:	ΑE,	AG,	AL,	AM,	AT,	AU,	AZ,	BA	, BB,	BG,	BR,	BY,	BZ,	CA,	CH,	CN,	
										, EC,								
		GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP	, KE,	KG,	KP,	KR,	KZ,	LC,	LK,	LR,	
		LS,	LT,	LU,	LV,	MA,	MD,	MG,	MK	, MN,	MW,	MX,	MZ,	NO,	NZ,	PL,	PT,	
		RO,	RU,	SD,	SE,	SG,	SI,	SK,	SL	, TJ,	TM,	TR,	TT,	ΤZ,	UA,	UG,	US,	
		UZ,	VN,	YU,	ZA,	ZW,	AM,	AZ,	BY	, KG,	ΚZ,	MD,	RU,	ТJ,	TM			
	RW	: GH,	GM,	KE,	LS,	MW,	ΜZ,	SD,	SL	, SZ,	TZ,	UG,	ZW,	ΑT,	BE,	CH,	CY,	
		DE,	DK,	ES,	FI,	FR,	GB,	GR,	ΙE	, IT,	LU,	MC,	NL,	PT,	SE,	TR,	BF,	
		ВJ,	CF,	CG,	CI,	CM,	GΑ,	GN,	GW	, ML,	MR,	ΝE,	SN,	TD,	TG			
AU 2001068958 A5 2002				2002	0114	AU 2001-68958						20010704 .						
E:	P 130	1187		A	2	2003	0416			EP 20	01-9	4721	1	2001	0704			
	R:	ΑT,	BE,	CH,	DE,	DK,	ES,	FR,	GB	, GR,	IT,	LI,	LU,	NL,	SE,	MC,	PT,	
		ΙE,	SI,	LT,	LV,	FI,	RO,	MK,	CY	, AL,	TR							
Ų:	S 200	21610	01	Α										2001				
PRIORI'	TY AP	PLN.	INFO	. :					DK :	2000-	1040		Α	2000	0704			
									US :	2000-	2232	40P	P	2000	0704			
								,	WO :	2001-	DK46	7	W	2001	0704			
OTHER :	SOURC	E(S):			MAR	PAT	136:	1024	02									

OTHER SOURCE(S): MA

MARPAT 136:102402

GI

AB Title compds. [I; m, n = 1, 2; Q = (CH2)n; Q1 = (CH2)m; R1 = CO, CS, (substituted) alkyl, alkenyl, alkynyl, cycloalkyl, cycloheteroalkyl, aryl, aralkyl, heteroaryl, heteroarylalkyl; R2 = H, (substituted) alkyl, alkenyl, alkynyl, cycloalkyl, cycloheteroalkyl, aryl, aralkyl, heteroarylalkyl, heteroaryl, cyano, halo, OH, NO2, SH, SR5, SOR5, SO2R5,

CO2H, CO2R4, CON(R5)2, alkoxy, alkenyloxy, alkynyloxy, aryloxy, heteroaryloxy; R3 = H, alkyl, alkenyl, alkynyl, cycloalkyl, cycloheteroalkyl, aryl, aralkyl, heteroarylalkyl, heteroaryl, alkoxyalkyl, CO2H, cyano, NO2, halo, OH; R3R3C, R4R4C = spiro system; R4 = (substituted) alkyl, alkenyl, alkynyl, cycloalkyl, cycloheteroalkyl, aryl, heteroaryl, cyano, halo, OH, NO2, CF3, alkoxy, alkenyloxy, alkynyloxy, aryloxy, heteroaryloxy, etc.; R5 = H, (substituted) alkyl, alkenyl, alkynyl, cycloalkyl, cycloheteroalkyl, aryl, arylalkyl, heteroaryl, heteroarylalkyl; R6 = H, (substituted) alkyl, alkenyl, alkynyl, cycloalkyl, cycloheteroalkyl, aryl, heteroaryl; R7 = H, (substituted) alkyl, alkenyl, alkynyl, cycloalkyl, cycloheteroalkyl, aryl, heteroaryl; R9 = H, halo, (substituted) alkyl, aryl; R10 = H, halo; R9R10 = atoms to form a cyclopropyl ring], were prepd. as DPP-IV inhibitors (no data). Thus, 1,4-diazepan-6-ylmethanol, 7-benzyl-8-chloro-1,3-dimethyl-3,7dihydropurine-2,6-dione, and K2CO3 were stirred at room temp. for 24 h, heated to 60.degree. for 3 h, heated to 95.degree. for 5 h, and heated to 120.degree. for 2 h to give 7-benzyl-8-(6-hydroxymethyl-1,4-diazepan-1-yl)-1,3-dimethyl-3,7-dihydropurine-2,6-dione.

IT 389060-65-5P 389062-41-3P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(prepn. of piperazinylpurinediones as inhibitors of dipeptidylpeptidase IV)

RN 389060-65-5 HCAPLUS

CN 1H-Purine-2,6-dione, 8-(hexahydro-1H-1,4-diazepin-1-yl)-3,7-dihydro-3-methyl-7-(phenylmethyl)-1-(3-phenyl-2-propenyl)- (9CI) (CA INDEX NAME)

$$\begin{array}{c|c} & \text{Me} \\ \hline \\ \text{O} & \text{N} & \text{N} \\ \hline \\ \text{Ph-CH} = \text{CH-CH}_2 & \text{O} \\ \hline \\ \text{O} & \text{CH}_2 - \text{Ph} \\ \end{array}$$

RN 389062-41-3 HCAPLUS

CN 1H-Purine-2,6-dione, 8-(hexahydro-1H-1,4-diazepin-1-yl)-3,7-dihydro-3-methyl-7-(phenylmethyl)-1-(3-phenyl-2-propenyl)-, trifluoroacetate (9CI) (CA INDEX NAME)

CM 1

CRN 389060-65-5 CMF C27 H30 N6 O2

$$\begin{array}{c|c} \text{Me} & & \\ \text{N} & & \\ \text{N} & & \\ \text{N} & & \\ \text{N} & & \\ \text{Ph-CH-CH-CH2} & & \\ \text{O} & & \\ \text{CH2-Ph} & \\ \end{array}$$

CM 2

CRN 76-05-1 CMF C2 H F3 O2

=> d 16 1-10

ANSWER 1 OF 10 REGISTRY COPYRIGHT 2003 ACS L6

RN 477333-67-8 REGISTRY

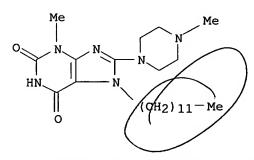
1H-Purine-2,6-dione, 7-dodecyl-3,7-dihydro-3-methyl-8-(4-methyl-1-CN

piperazinyl) - (9CI) (CA INDEX NAME)

FS 3D CONCORD

C23 H40 N6 O2 MF

SR CAS Registry Services



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

ANSWER 2 OF 10 REGISTRY COPYRIGHT 2003 ACS L6

454706-71-9 REGISTRY

1H-Purine-2,6-dione, 7-(2,3-dimethyl-2-butenyl)-3,7-dihydro-1,3-dimethyl-8-

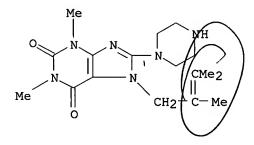
(1-piperazinyl) - (9CI) (CA INDEX NAME)

3D CONCORD

MF C17 H26 N6 O2

SR CA

LC STN Files: CA, CAPLUS



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

1 REFERENCES IN FILE CA (1957 TO DATE)

1 REFERENCES IN FILE CAPLUS (1957 TO DATE)

L6 ANSWER 3 OF 10 REGISTRY COPYRIGHT 2003 ACS

RN 402470-56-8 REGISTRY

CN 1H-Purine-2,6-dione, 3,7-dihydro-8-[4-(2-hydroxyethyl)-1-piperazinyl]-3methyl-7-nonyl- (9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C21 H36 N6 O3 SR Chemical Library

LC STN Files: CHEMCATS

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L6 ANSWER 4 OF 10 REGISTRY COPYRIGHT 2003 ACS

RN · 378201-33-3 REGISTRY

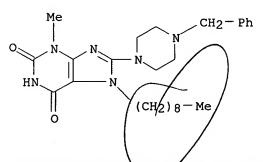
CN 1H-Purine-2,6-dione, 3,7-dihydro-3-methyl-7-nonyl-8-[4-(phenylmethyl)-1-piperazinyl]- (9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C26 H38 N6 O2

SR Chemical Library

LC STN Files: CHEMCATS



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

- L6 ANSWER 5 OF 10 REGISTRY COPYRIGHT 2003 ACS
- RN 377060-82-7 REGISTRY
- CN 1H-Purine-2,6-dione, 3,7-dihydro-3-methyl-7-octyl-8-[4-(phenylmethyl)-1-piperazinyl]- (9CI) (CA INDEX NAME)
- FS 3D CONCORD
- MF C25 H36 N6 O2
- SR Chemical Library
- LC STN Files: CHEMCATS

## \*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

- L6 ANSWER 6 OF 10 REGISTRY COPYRIGHT 2003 ACS
- RN 371927-37-6 REGISTRY
- CN 1H-Purine-2,6-dione, 7-hexyl-3,7-dihydro-3-methyl-8-(1-piperazinyl)- (9CI) (CA INDEX NAME)
- FS 3D CONCORD
- MF C16 H26 N6 O2
- SR Chemical Library
- LC STN Files: CHEMCATS

# \*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

- L6 ANSWER 7 OF 10 REGISTRY COPYRIGHT 2003 ACS
- RN 329702-29-6 REGISTRY
- CN 1H-Purine-2,6-dione, 3,7-dihydro-3-methyl-7-nonyl-8-(1-piperazinyl)- (9CI) (CA INDEX NAME)
- FS 3D CONCORD
- MF C19 H32 N6 O2
- SR Chemical Library
- LC STN Files: CHEMCATS

## \*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L6 ANSWER 8 OF 10 REGISTRY COPYRIGHT 2003 ACS

RN 313471-56-6 REGISTRY

CN 1H-Purine-2,6-dione, 7-heptyl-3,7-dihydro-3-methyl-8-(4-methyl-1-piperazinyl)- (9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C18 H30 N6 O2

SR Chemical Library

LC STN Files: CHEMCATS

$$\begin{array}{c|c} Me & Me \\ \hline N & N & N \\ \hline N & N & N \\ \hline O & (CH_2)_6-Me \end{array}$$

### \*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L6 ANSWER 9 OF 10 REGISTRY COPYRIGHT 2003 ACS

RN 304876-71-9 REGISTRY

CN 1H-Purine-2,6-dione, 7-hexadecyl-3,7-dihydro-3-methyl-8-(1-piperazinyl)-(9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C26 H46 N6 O2

SR Chemical Library

LC STN Files: CHEMCATS

$$\begin{array}{c|c} Me & & NH \\ \hline N & N & NH \\ \hline N & NH \\ N & NH \\ \hline N & NH \\ N & NH \\ \hline N & NH \\ N & NH \\ \hline N & NH \\ N & NH \\ \hline N & NH \\ N & NH \\ \hline N & NH \\ N & NH \\ \hline N & NH \\ N & NH \\ \hline N & NH \\ N & NH \\ \hline N & NH \\ N & NH \\ \hline N$$

### \*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L6 ANSWER 10 OF 10 REGISTRY COPYRIGHT 2003 ACS

RN 106939-21-3 REGISTRY

CN 1H-Purine-2,6-dione, 7-heptyl-3,7-dihydro-3-methyl-8-(1-piperazinyl)-(9CI) (CA INDEX NAME)

FS 3D CONCORD

# Berch 09/935,149

MF C17 H28 N6 O2

SR CA

LC STN Files: CA, CAPLUS, CASREACT, CHEMCATS, TOXCENTER

# \*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

- 1 REFERENCES IN FILE CA (1957 TO DATE)
- 1 REFERENCES IN FILE CAPLUS (1957 TO DATE)